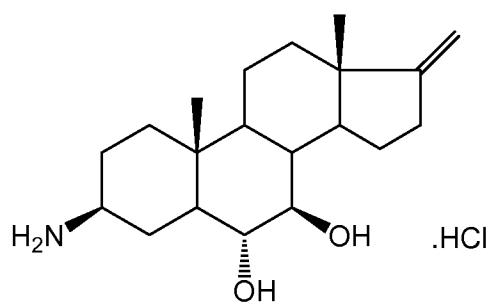


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1) (Original) Crystalline form A of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride corresponding to the structure:



characterized by the fact that the indexing of the first 30 lines of the powder X-ray diffraction pattern diagram at 295 K is:

| <i>h</i> | <i>k</i> | <i>l</i> | lattice spacing (Å) | 2 theta "mean λ Cu $\text{K}\alpha$ " 1.54184 Å |
|----------|----------|----------|------------------------|---|
| 1 | 0 | 0 | 16.058 | 5.50 |
| 0 | 0 | 1 | 9.011 | 9.82 |
| 2 | 0 | 0 | 8.029 | 11.02 |
| -1 | 0 | 1 | 7.872 | 11.24 |
| 1 | 0 | 1 | 7.844 | 11.28 |
| 1 | 1 | 0 | 6.413 | 13.81 |
| -2 | 0 | 1 | 6.007 | 14.75 |
| 2 | 0 | 1 | 5.982 | 14.81 |

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| h | k | l | lattice spacing (\AA) | 2 theta "mean" λCu $\text{K}\alpha$ " 1.54184 \AA |
|-----|-----|-----|-------------------------------------|---|
| 0 | 1 | 1 | 5.526 | 16.04 |
| 3 | 0 | 0 | 5.353 | 16.56 |
| 2 | 1 | 0 | 5.274 | 16.81 |
| -1 | 1 | 1 | 5.229 | 16.96 |
| 1 | 1 | 1 | 5.221 | 16.98 |
| -3 | 0 | 1 | 4.610 | 19.25 |
| 3 | 0 | 1 | 4.594 | 19.32 |
| -2 | 1 | 1 | 4.557 | 19.48 |
| 2 | 1 | 1 | 4.546 | 19.53 |
| 0 | 0 | 2 | 4.506 | 19.70 |
| -1 | 0 | 2 | 4.343 | 20.45 |
| 1 | 0 | 2 | 4.333 | 20.50 |
| 3 | 1 | 0 | 4.251 | 20.90 |
| 4 | 0 | 0 | 4.014 | 22.14 |
| -2 | 0 | 2 | 3.936 | 22.59 |
| 2 | 0 | 2 | 3.922 | 22.67 |
| -3 | 1 | 1 | 3.850 | 23.11 |
| 3 | 1 | 1 | 3.840 | 23.17 |
| 0 | 1 | 2 | 3.788 | 23.49 |
| -1 | 1 | 2 | 3.690 | 24.12 |
| 1 | 1 | 2 | 3.684 | 24.16 |
| -4 | 0 | 1 | 3.673 | 24.23 |

2) (Original) Crystalline form A of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride characterized by the fact that the unit cell is monoclinic (space group P2, Z=2) and the unit cell parameters at T = 295 K are:

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$$\begin{array}{ll}
 a = 16.058(2) \text{ \AA}, & \beta = 90.24(2)^\circ \\
 b = 6.995(1) \text{ \AA}, & V = 1012.2 \text{ \AA}^3 \\
 c = 9.011(2) \text{ \AA} & \text{density} = 1.168
 \end{array}$$

3) (Original) Di-hydrated crystalline form B of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride, characterized by the fact that the indexing of the lines of the powder X-ray diffraction pattern diagram at 295 K is:

| h | k | l | Lattice spacing (\AA) | 2 theta "mean λ Cu $K\alpha$ " 1.54184 \AA |
|-----|-----|-----|-------------------------------------|---|
| 0 | 1 | 0 | 17.770 | 4.97 |
| 0 | 2 | 0 | 8.885 | 9.96 |
| 1 | 0 | 0 | 8.667 | 10.21 |
| 1 | 1 | 1 | 8.509 | 10.40 |
| -1 | 1 | 0 | 7.227 | 12.25 |
| 1 | 2 | 0 | 6.960 | 12.72 |
| 0 | 0 | 1 | 6.778 | 13.06 |
| 0 | -1 | 1 | 6.777 | 13.06 |
| 0 | 1 | 1 | 5.966 | 14.85 |
| 0 | -2 | 1 | 5.964 | 14.85 |
| 0 | 3 | 0 | 5.923 | 14.96 |
| -1 | 2 | 0 | 5.651 | 15.68 |
| -1 | -1 | 1 | 5.446 | 16.28 |
| 1 | 0 | 1 | 5.441 | 16.29 |
| 1 | 3 | 0 | 5.438 | 16.30 |
| 1 | 0 | 1 | 5.243 | 16.91 |
| 1 | -1 | 1 | 5.238 | 16.93 |
| -1 | -2 | 1 | 5.172 | 17.15 |
| 1 | 1 | 1 | 5.168 | 17.16 |
| 0 | 2 | 1 | 4.953 | 17.91 |
| 0 | -3 | 1 | 4.952 | 17.91 |

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| h | k | l | Lattice spacing (\AA) | 2 theta "mean $\lambda\text{Cu K}\alpha$ " 1.54184 \AA |
|----|----|---|-------------------------------------|---|
| -1 | 1 | 1 | 4.695 | 18.90 |
| 1 | -2 | 1 | 4.690 | 18.92 |
| -1 | -3 | 1 | 4.594 | 19.32 |
| 1 | 2 | 1 | 4.591 | 19.33 |
| -1 | 3 | 0 | 4.481 | 19.82 |
| 0 | 4 | 0 | 4.443 | 19.99 |
| 2 | 1 | 0 | 4.425 | 20.07 |
| 2 | 0 | 0 | 4.334 | 20.49 |
| 1 | 4 | 0 | 4.331 | 20.51 |

4) (Original) Di-hydrated crystalline form B of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride, characterized by the fact that the unit cell is triclinic (space group P1, Z=1) and the unit cell parameters at T = 295 are:

$$\begin{aligned}
 a &= 8.856(2) \text{ \AA}, & \alpha &= 100.76(1)^\circ \\
 b &= 18.482(1) \text{ \AA}, & \beta &= 90.06(1)^\circ \\
 c &= 6.904(2) \text{ \AA}, & \gamma &= 78.35(l) \\
 & & V &= \text{do } 1086.5 \text{ \AA}^3 \\
 & & \text{density} &= 1.198
 \end{aligned}$$

5) (Original) Monohydrate crystalline form C of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride, characterized by the fact that the indexing of the first 30 lines of the powder X-ray diffraction pattern diagrams at 295K is :

| h | k | l | Lattice spacing (\AA) | 2 theta "mean $\lambda\text{Cu K}\alpha$ " 1.54184 \AA |
|---|---|---|-------------------------------------|---|
| 0 | 1 | 0 | 20.875 | 4.23 |

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| h | k | l | Lattice spacing (\AA) | 2 theta "mean $\lambda_{\text{Cu K}\alpha}$ " 1.54184 \AA |
|-----|-----|-----|-------------------------------------|--|
| 0 | 2 | 0 | 10.437 | 8.47 |
| 1 | 0 | 0 | 7.049 | 12.56 |
| 0 | 3 | 0 | 6.958 | 12.72 |
| 0 | 0 | 1 | 6.922 | 12.79 |
| 0 | -1 | 1 | 6.845 | 12.93 |
| -1 | 1 | 0 | 6.780 | 13.06 |
| 1 | 1 | 0 | 6.581 | 13.46 |
| 0 | 1 | 1 | 6.325 | 14.00 |
| 0 | -2 | 1 | 6.155 | 14.39 |
| -1 | 2 | 0 | 5.980 | 14.81 |
| 1 | 2 | 0 | 5.712 | 15.51 |
| -1 | 0 | 1 | 5.604 | 15.81 |
| -1 | -1 | 1 | 5.506 | 16.10 |
| 0 | 2 | 1 | 5.447 | 16.27 |
| -1 | 1 | 1 | 5.323 | 16.66 |
| 0 | -3 | 1 | 5.267 | 16.83 |
| 0 | 4 | 0 | 5.219 | 16.99 |
| -1 | -2 | 1 | 5.083 | 17.45 |
| -1 | 3 | 0 | 5.079 | 17.46 |
| 1 | 3 | 0 | 4.834 | 18.35 |
| -1 | 2 | 1 | 4.804 | 18.47 |
| 0 | 3 | 1 | 4.612 | 19.24 |
| -1 | -3 | 1 | 4.516 | 19.66 |
| 1 | -1 | 1 | 4.474 | 19.84 |
| 1 | 0 | 1 | 4.465 | 19.88 |
| 0 | -4 | 1 | 4.459 | 19.91 |
| -1 | 4 | 0 | 4.297 | 20.67 |
| 1 | -2 | 1 | 4.290 | 20.71 |

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| h | k | l | Lattice spacing (\AA) | 2 theta "mean $\lambda\text{Cu K}\alpha$ " 1.54184 \AA |
|---|---|---|-------------------------------------|---|
| 1 | 1 | 1 | 4.266 | 20.82 |

6) (Original) Monohydrate crystalline form C of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride, characterized by the fact that the unit cell is triclinic (space group P1, Z=1) and the unit cell parameters T = 295 K are:

$$\begin{aligned}
 a &= 7.2328(5) \text{ \AA} & \alpha &= 97.135(6)^\circ \\
 b &= 21.063 (2) \text{ \AA} & \beta &= 102.653(5)^\circ \\
 c &= 7.1563(5) \text{ \AA} & \gamma &= 91.177(6)^\circ \\
 & & V &= 1054.2 \text{ \AA}^3 \\
 & & \text{density} &= 1.178
 \end{aligned}$$

7) (Original) A process for the preparation of form A as defined in Claims 1 or 2, characterized by the fact that crystallization takes place in a mixture of alcohol and ether and particularly in an isopropylic methanol-ether mixture.

8) (Original) A process for the preparation of form C as defined in Claim 5 or 6, characterized by the fact that 250 mg of compound of formula (1) are dissolved at ambient temperature in a solvent such as methyl ethyl ketone (MEK); and then transferred in water by azeotropic distillation at constant volume and equilibration at a relative humidity above 97%.

9) (Currently Amended) As medications, crystalline forms A, B or C as defined by
 | Claims 1 to 86.

10) (Currently Amended) A pharmaceutical composition characterized by the fact that it comprises form A of 3-beta-amino-17-methylene-androstane-6-alpha,7-beta-diol hydrochloride in a pure state or possibly in combination with either one of/or both

crystalline forms B or C and/or in combination with any compatible and pharmaceutically acceptable ~~additive~~ excipient or inert diluent.

11) (Currently Amended) ~~Application of the crystalline forms as defined by one any of the Claims 1 to 8 for the preparation of a medicament for the treatment of inflammatory diseases~~ A method of treating inflammatory diseases, wherein the method comprises administering a pharmaceutical composition comprising one or more crystalline forms A, B or C as defined by Claims 1 to 6 to a human.